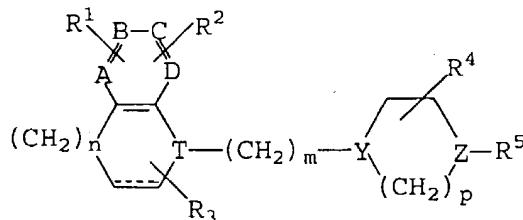


AMENDMENTS TO THE CLAIMS

1. (Currently Amended) A ~~1,4-substituted cyclic amine~~  
~~derivative represented by the following~~ compound of formula (I):



(I)

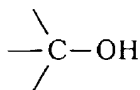
wherein A, B, C, D, and T are the same or different from one another and each represents methine or nitrogen, provided that one and only one of them represents nitrogen;

the bond represented by the following formula:

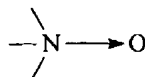


represents a single or double bond;

Y and Z are the same or different from each other and each represents methine, nitrogen, a group represented by the following formula:



or a group represented by the following formula:



provided at least one of them represents nitrogen;

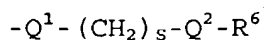
R<sup>1</sup> and R<sup>2</sup> are the same or different from each other and each represents hydrogen, halogeno, hydroxy, lower alkylsulfonylaminoalkyl, lower halogenated-alkylsulfonylaminoalkyl, 2-pyrrolidinon-1-yl, 1-hydroxy-1-(methoxypyridyl)methyl, methoxypyridylcarbonyl, 1,3-propanesultum-2-yl, lower hydroxypiperidyl-carbonylalkyl, lower hydroxyalkylamidoalkyl, lower halogenated-alkylamidoalkyl, lower dihalogenatedalkylamidoalkyl, ~~lower—heteroaryl-amidoalkyl,~~ lower hydroxyalkylamidoalkyl, optionally substituted amino, nitro, lower alkyl, lower alkoxy, lower acyl, lower alkoxyalkoxy, cyano, lower alkylsulfonyl, sulfonylamido, hydroxy-lower alkyl, hydroxy-lower alkoxy, lower alkoxycarbonylamino, lower alkylsulfonylamino, N-lower alkylalkylsulfonylamino, lower acylamino, optionally substituted aminoalkyl, optionally N-substituted lower acylaminoalkyl, optionally substituted aryl, optionally substituted arylsulfonylamino, lower alkylsulfonyloxy, hydroxyiminomethyl, (2-pyrrolidon1-yl)methyl, (2-piperidon-1-yl)methyl, optionally substituted heteroaryl, optionally substituted aralkyl, optionally substituted heteroarylalkyl, cycloalkylcarbonylaminoalkyl, optionally substituted ureido, optionally substituted ureido-lower alkyl, succinimido, (succinimido-1-yl)-lower alkyl, amido, optionally substituted carbamoyl, optionally substituted carbamoyl-lower alkyl, optionally substituted

thiocarbamoyl lower alkyl, formyl, aromatic acyl, heteroarylcarbonyl, halogenated lower alkyl, (2-imidazolidinon-1-yl)methyl, (2,4-imidazolidinedion-3-yl)methyl, (2-oxazolidon-3-yl)methyl, (glutarimido-1-yl)methyl, optionally substituted heteroarylhydroxyalkyl, cyano-lower alkyl, 1-hydroxy lower cycloalkyl, (2,4-thiazolidinedion-3-yl)methyl, optionally substituted 4-piperidylmethyl, heteroarylacyl, pyrrolidinylcarbonyl-lower alkyl, optionally substituted aminosulfonylalkyl, carboxy-lower alkyl, or lower alkylamidoalkyl; or alternatively  $R^1$  and  $R^2$  together may form optionally substituted ~~alicyclic~~, optionally substituted heterocycle or alkylenedioxy, provided these rings may be substituted;

$R^3$  represents hydrogen, halogeno, lower alkyl, hydroxy, hydroxy-lower alkyl, lower alkoxy, formyl, optionally substituted aralkyloxy, hydroxy-lower alkoxy, optionally substituted sulfamoyl, or optionally N-substituted sulfamoyl-lower alkyl;

$R^4$  represents hydrogen, lower alkyl, hydroxy-lower alkyl, lower alkoxyalkyl, optionally aryl-substituted aryloxyalkyl, or optionally aryl-substituted aralkyloxyalkyl;

$R^5$  represents lower alkyl, lower acyl, lower alkoxy carbonyl, aromatic acyl, or a group represented by the following formula:



wherein  $Q^1$  and  $Q^2$  are both single bonds, or one of them is a single bond while the other represents oxygen, carbonyl, a group represented by  $-NHCO-$ , a group represented by  $-NHSO_2-$ , or a group represented by  $>CH-R^7$ , wherein  $R^7$  represents hydroxy, lower alkyl or halogeno:

s represents 0 or an integer of 1 to 6; and

$R^6$  represents optionally substituted aryl, ~~optionally substituted heteroaryl~~, optionally substituted benzoheteroaryl, 1,4-benzodioxanyl, 1,3-benzodioxolyl, benzothiazolyl, or cyano;

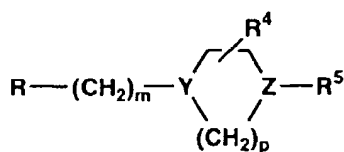
n represents 1;

m represents 0 or an integer of 1 to 6; and

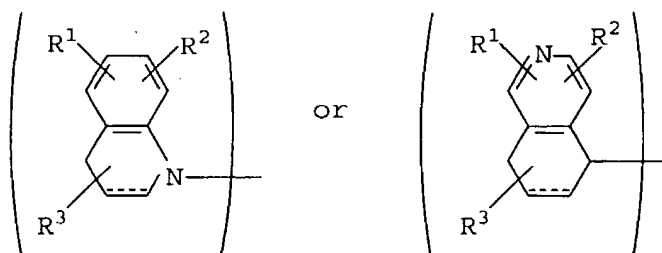
p represents an integer of 1 to 3,

or a and pharmacologically acceptable salt ~~salts~~ thereof.

2. (Currently Amended) A ~~1,4-substituted cyclic amine derivative represented by the following~~ compound of formula:



wherein R represents a substituent of the formula:



wherein the bond represented by the following formula:



and  $R^1$ ,  $R^2$ ,  $R^3$ ,  $R^4$ ,  $R^5$ , Y, Z, m, and p are each as defined in claim 1, and pharmacologically acceptable salts thereof.

3. (Currently Amended) The ~~1,4-substituted cyclic amine derivative as set forth in~~ compound of claim 1 or a pharmacologically acceptable salt thereof, wherein m is 0 and p is 2.

4. (Currently Amended) The ~~1,4-substituted cyclic amine derivative as set forth in~~ compound of claim 1 or a pharmacologically acceptable salt thereof, wherein Y is methine and Z is nitrogen.

5. (Currently Amended) The ~~1,4-substituted cyclic amine derivative as set forth in~~ compound of claim 1 or a pharmacologically acceptable salt thereof, which is a compound selected from among the following ones:

(267) 1-{1-[2-(4-methoxyphenyl)ethyl]piperidin-4-yl}-7-methoxy-1,2,3,4-tetrahydroquinoline,

(268) 1-{1-[2-(4-fluorophenyl)ethyl]piperidin-4-yl}-7-methoxy-1,2,3,4-tetrahydroquinoline,

(269) 1-[1-(4-cyanopropyl)piperidin-4-yl]-7-methoxy-1,2,3,4-tetrahydroquinoline,

(270) 1-{1-[2-(2-thienyl)ethyl]piperidin-4-yl}-7-methoxy-1,2,3,4-tetrahydroquinoline,

(271) 1-{1-[2-(4-fluorophenyl)ethyl]piperidin-4-yl}-7,8-dimethoxy-1,2,3,4-tetrahydroquinoline,

(272) 1-{1-[2-(4-fluorophenyl)ethyl]piperidin-4-yl}-7,8-methylenedioxy-1,2,3,4-tetrahydroquinoline,

(273) 1-{1-[2-(4-fluorophenyl)ethyl]piperidin-4-yl}-7-methoxy-8-methyl-1,2,3,4-tetrahydroquinoline,

(274) 1-{1-[2-(4-fluorophenyl)-2-oxoethyl]piperidin-4-yl}-7-methoxy-1,2,3,4-

tetrahydroquinoline,

(275) 1-{1-[2-(4-fluorophenyl)-2-hydroxyethyl]piperidin-4-yl}-7-methoxy-1,2,3,4-

tetrahydroquinoline,

(276) 1-{1-[2-(4-fluorophenyl)-2-fluoroethyl]piperidin-4-yl}-7-methoxy-1,2,3,4-

tetrahydroquinoline, and

(283) 5-{4-[2-(4-fluorophenyl)ethyl]piperazin-1-yl}-5,6,7,8-tetrahydroisoquinoline.

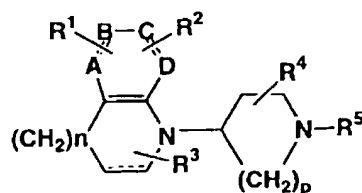
6. (Currently Amended) A pharmaceutical composition comprising a therapeutically effective amount of the ~~1,4-substituted cyclic amine derivative or salt as set forth in~~ compound of claim 1 or its salt in combination with a pharmaceutically acceptable carrier.

7. (Currently Amended) An agent for treating, and ~~ameliorating, and preventing~~ diseases against which serotonin antagonism is efficacious, which ~~contain~~ contains as the active ingredient the ~~1,4-substituted cyclic amine derivative as set forth in~~ compound of claim 1 or a pharmacologically acceptable salt thereof.

8. (Currently Amended) A compound ~~An agent~~ for treating, and ameliorating, ~~and preventing~~ spastic paralysis, which comprises ~~contain~~ as the active ingredient the ~~1,4-substituted cyclic amine derivative as set forth in~~ an effective amount of the compound of claim 1 or a pharmacologically acceptable salt thereof.

9. (Currently Amended) A muscle relaxant composition which contains as the active ingredient the ~~1,4-substituted cyclic amine derivative as set forth in~~ an effective amount of the compound of claim 1 or a pharmacologically acceptable salt thereof.

10. (Currently Amended) A process for producing a ~~1,4-substituted cyclic amine derivative represented by the following~~ compound of the formula:



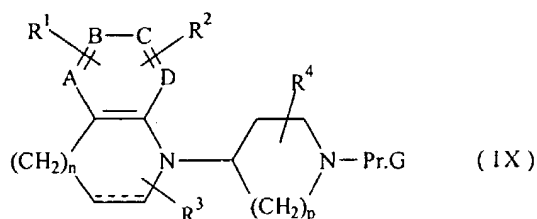
wherein the bond represented by the following formula:



and A, B, C, D, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, n, and p are each as defined in claim 1, which comprises removing, if necessary, the



protecting group from a 1,4-substituted cyclic amine derivative  
(IX) represented by the following formula:

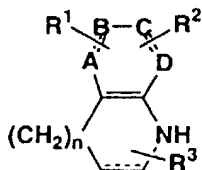


wherein the bond represented by the following formula:



and A, B, C, D, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, n, and p are each as defined in claim 1; and Pr.G represents hydrogen or a protecting group, and then reacting the same with L-R<sup>5</sup> wherein R<sup>5</sup> is as defined in claim 1; and L represents a leaving group.

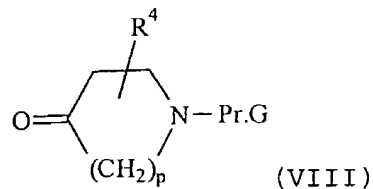
11. (Currently Amended) A process for producing 1,4-substituted cyclic amine derivative (X), as set forth in a compound of claim 1, which comprises reacting a fused cyclic amine represented by the following formula:



wherein the bond represented by the following formula:

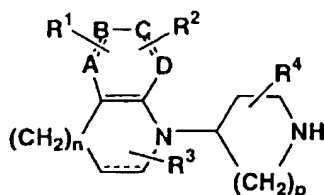


and A, B, C, D, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup> and n are each as defined in claim 1  
with a cyclic ketone (VIII) represented by the following  
formula:



wherein R<sup>4</sup>, p, and Pr.G are each as defined in claim 1  
in the presence of a reducing agent to thereby give a 1,4-  
substituted cyclic amine derivative (IX), removing, if necessary,  
the protecting group therefrom and further reacting the same with  
L-R<sup>5</sup>.

12. (Currently Amended) A ~~4-substituted cyclic amine~~  
~~derivative represented by the following~~ compound of the formula:



wherein the bond represented by the following formula:



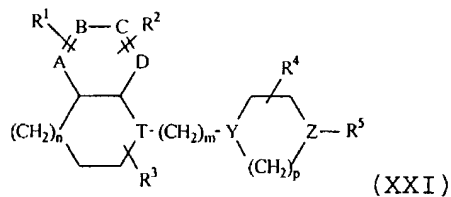
and A, B, C, D, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, n, and p are each as defined in claim 1, provided that the case where R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> are all hydrogen atoms is excluded.

13. (Currently Amended) A method for ~~treating a disease to which antagonizing serotonin antagonism is efficacious~~, which comprises administering an effective dose of the ~~1,4-substituted cyclic amine derivative as set forth in~~ compound of claim 1, or a pharmacologically acceptable salt thereof, to a person in need of such treatment.

14. (Currently Amended) The ~~1,4-substituted cyclic amine derivative as set forth in~~ compound of claim 1, in which the bond represented by the following formula in the formula (I):

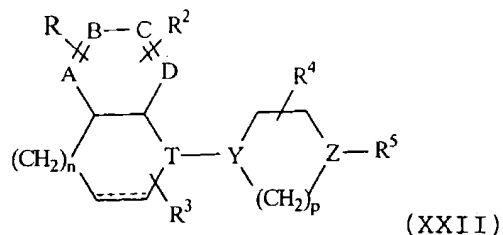
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is a single bond, represented by the formula (XXI):



or a pharmacologically acceptable salt thereof.

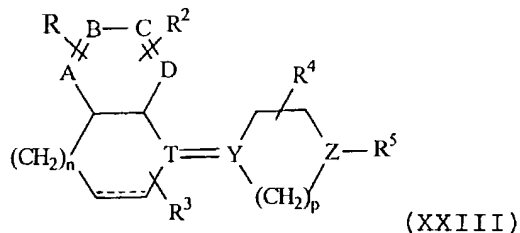
15. (Currently Amended) The ~~1,4-substituted cyclic amine~~ derivative ~~as set forth in~~ compound of claim 1, in which m is 0 in the formula (I), represented by the formula (XXII):



or a pharmacologically acceptable salt thereof.

16. (Currently Amended) The ~~1,4-substituted cyclic amine~~ derivative ~~as set forth in~~ compound of Claim 1, in which m is 1 to 6 in the formula (I) or a pharmacologically acceptable salt thereof.

17. (Currently Amended) A ~~1,4-substituted cyclic amine~~ derivative ~~represented by~~ compound of the formula (XXIII):



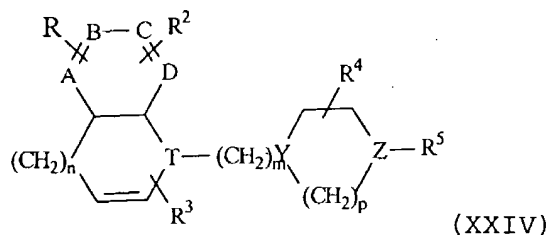
or a pharmacologically acceptable salt thereof.

18. (Currently Amended) The ~~1,4-substituted cyclic amine~~

~~derivative as set forth in~~ compound of claim 1, in which the bond represented by the following formula in the formula (I):

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is a double bond, represented by the formula (XXIV):



or a pharmacologically acceptable salt thereof.

19. (Currently Amended) The ~~1,4-substituted cyclic amine~~ derivative as set forth in compound of claim 1, in which the T is nitrogen.